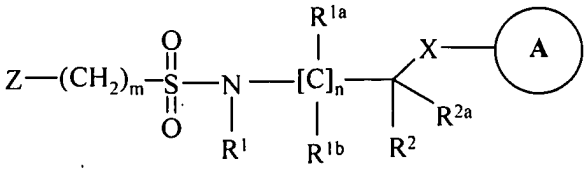
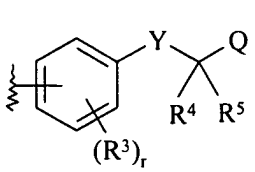
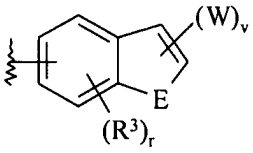
Amendments to the Claims

1. (Currently Amended) A compound having a structural Formula I,

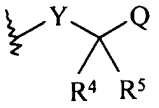


I

or ~~and~~ pharmaceutically acceptable salts~~, solvates, hydrates~~ or stereoisomers thereof, wherein:

A is:  or  ;

E is selected from the group consisting of~~:~~ O, S and~~or~~ NR^{14};

W is selected from the group consisting of~~:~~  , hydrogen, C_1-C_6 alkyl, $(CH_2)_n$-C_3-C_6 cycloalkyl, haloalkyl and~~or~~ acyl;

Q is: $-C(O)OR^6$ or R^{6A};

X is selected from the group consisting of~~:~~ a bond, CH_2, O, S and~~or~~ $S[O]_p$;

Y is selected from the group consisting of~~:~~ a bond, S, CH_2 and~~or~~ O;

Z is: benzothiophene; ~~a) aliphatic group,~~

~~b) aryl,~~

~~c) a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S,~~

~~d) bi-aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl,~~

~~e) bi-heteroaryl, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and~~

~~f) heterocyclyl;~~

wherein the benzothiophene~~aliphatic group, aryl, heteroaryl, bi-aryl, bi-heteroaryl and heterocyclyl being~~ is optionally substituted with one or more groups independently selected from R^{15};

m and n' are each independently selected from the group consisting of~~:~~ 0, 1, 2, 3 and~~or~~ 4;

n is selected from the group consisting of~~:~~ 0, 1, 2 and~~or~~ 3;

p is: 1 or 2;

r is selected from the group consisting of~~:~~ 1, 2, 3 and~~or~~ 4;

v is: 1 or 2;

R^1 is selected from the group consisting of~~:~~ hydrogen, ~~wherein when Z is phenyl or naphthyl and R^2 is H, R^1 is not H,~~

haloalkyl,

C_1-C_6 alkyl,

C_1-C_6 alkyl-C_1-C_6 alkoxy,

C_1-C_6 alkyl-aryl,

C_2-C_6 alkenyl,

C_2-C_6 alkynyl,

$(CH_2)_n$-C_3-C_6 cycloalkyl,

C_1-C_6 alkoxy,

aryl, and~~or~~

R^1 and R^2 together being a 5- to 8-membered heterocyclyl ring;~~, and~~

wherein alkyl, aryl, alkenyl, alkynyl, cycloalkyl and alkoxy being optionally substituted with one or more groups independently selected from R^{15};

R^{1a} and R^{1b} are each independently selected from the group consisting of~~:~~

hydrogen,

C_1-C_6 alkyl, and~~or~~

R^1 and R^{1a}, R^1 and R^{1b}, R^2 and R^{1a}, R^2 and R^{1b} or R^{1a} and R^{1b} together being a 3- to 6-membered heterocyclyl or carbocyclyl ring wherein~~where~~ at least one of R^{1a} and R^{1b} is not hydrogen;

R^2 is selected from the group consisting of —hydrogen,

haloalkyl,

C_1-C_6 alkyl,

C_1-C_6 alkyl-C_1-C_6 alkoxy,

C_1-C_6 alkyl-aryl,

C_2-C_6 alkenyl,

C_2-C_6 alkynyl,

$(CH_2)_n$-C_3-C_6 cycloalkyl,

C_1-C_6 alkoxy,

aryl, and~~or~~

R^1 and R^2 together being a 5- to 8-membered heterocyclyl ring;~~, and~~

wherein alkyl, aryl, alkenyl, alkynyl, cycloalkyl and alkoxy being optionally substituted with one or more groups independently selected from R^{15};

R^{2a} is selected from the group consisting of~~:~~ hydrogen, halo,~~ or~~ C_1-C_6 alkyl and wherein R^2 and R^{2a} together being a 3- to 8-membered ring;~~ and~~ wherein alkyl being optionally substituted with one or more groups independently selected from R^{15};

R^3 is selected from the group consisting of~~:~~ — hydrogen,

halo,

cyano,

haloalkyl,

C_1–C_6 alkyl,

$(CH_2)_n$-C_3-C_6 cycloalkyl,

(C_1–C_4 alkyl)-heterocyclyl,~~ wherein the heterocyclyl being optionally substituted with oxo,~~ and

(C_1–C_4 alkyl)-$NR^7C(O)_pR^9$;~~ and~~

wherein alkyl, cycloalkyl and heterocyclyl being optionally substituted with one or more groups independently selected from R^{15};

R^4 and R^5 are each independently selected from the group consisting of:

hydrogen,

halo,

C_1-C_6 alkyl

C_1-C_6 alkoxy;

aryloxy;

$N(R^8)_2$,

SR^8 and~~or~~

R^4 and R^5 together being a 3- to 8-membered ring;

R^6 isselected from the group consisting of~~:~~ hydrogen, C_1-C_6 alkyl and~~or~~ aminoalkyl;

R^{6A} is selected from the group consisting of~~:~~ carboxamide, C_1-C_3 alkylnitrile, sulfonamide, acylsulfonamide and~~or~~ tetrazole;

R^7 is selected from the group consisting of ~~:~~hydrogen and~~or~~ C_1-C_6 alkyl;

R^8 and R^9 are each independently selected from the group consisting of:

hydrogen, C_1-C_6 alkyl, aryl, heteroaryl, and~~or~~ heterocyclyl~~;, and~~

wherein aryl, heteroaryl and heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl and C_1-C_6 alkoxy;

R^{14} is selected from the group consisting of~~:~~ hydrogen, aryl, C_1-C_6 alkyl, and~~or~~ C_1-C_6 alkyl-$COOR^6$,~~and~~

wherein aryl and alkyl being optionally substituted with one or more groups independently selected from R^{15}; and

R^{15} is selected from the group consisting of~~:~~ hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl, C_1-C_6 alkoxy, $(CH_2)_n$-C_3-C_6 cycloalkyl, $N(R^8)_2$, $NR^8S(O)_2R^9$, $NR^8C(O)_pR^9$, $C(O)NR^8R^9$, $C(O)_pR^8$, SR^8, $S(O)_pR^8$ and~~or~~ $S(O)_2NR^8R^9$.

2. (Currently Amended) The compound Claim 1, wherein X and Y are respectively S and O; S and CH_2; or CH_2 and O.

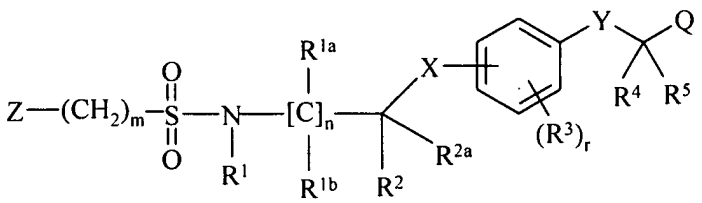
3. (Canceled)

4. (Canceled)

5. (Currently Amended) The compound of Claim 1~~Claim 4~~, wherein R^1 is C_3-C_6 alkyl or $(CH_2)_n$ C_3-C_6 cycloalkyl; R^2 and R^3 are each independently C_1-C_3 alkyl; and r is 1.

6. (Original) The compound Claim 5, wherein X is positioned para to Y; and R^3 is positioned ortho to Y.

7. (Currently Amended) A compound having a structural Formula II,



II

or~~and~~ pharmaceutically acceptable salts~~, solvates, hydrates~~ or stereoisomers thereof, wherein:

Q is: $-C(O)OR^6$ or R^{6A};

X is selected from the group consisting of: a bond, CH_2, O, S and~~or~~ $S[O]_p$;

Y is selected from the group consisting of: a bond, S, CH_2 and~~or~~ O;

Z is benzothiopene~~: a) aliphatic group. b) aryl, c) a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S,~~

~~d) bi-aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl.~~

~~e) bi-heteroaryl, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and~~

~~f) heterocyclyl;~~

wherein the benzothiophene~~aliphatic group, aryl, heteroaryl, bi-aryl, bi-heteroaryl and heterocyclyl being~~ is optionally substituted with one or more groups independently selected from R^{15};

m and n' are each independently selected from the group consisting of~~:~~ 0, 1, 2, 3 and~~or~~ 4;

n is selected from the group consisting of~~:~~ 0, 1, 2 and~~or~~ 3;

p is: 1 or 2;

r is selected from the group consisting of~~:~~ 1, 2, 3 and~~or~~ 4;

R^1 is selected from the group consisting of~~:~~ aryl,

haloalkyl,

C_1-C_6 alkyl,

C_1-C_6 alkyl-C_1-C_6 alkoxy,

C_1-C_6 alkyl-aryl,

C_2-C_6 alkenyl,

C_2-C_6 alkynyl,

$(CH_2)_n$-C_3-C_6 cycloalkyl,

C_1-C_6 alkoxy and~~or~~

R^1 and R^2 together being a 5- to 8-membered heterocyclyl ring~~; and~~

wherein alkyl, aryl, alkenyl, alkynyl, cycloalkyl and alkoxy being optionally substituted with one or more groups independently selected from R^{15};

R^{1a} and R^{1b} are each independently selected from the group consisting of~~:~~

hydrogen,

C_1-C_6 alkyl, and~~, or~~

R^1 and R^{1a}, R^1 and R^{1b}, R^2 and R^{1a}, R^2 and R^{1b} or~~or~~ R^{1a} and R^{1b} together being a 3- to 6-membered heterocyclyl or carbocyclyl ring where at least one of R^{1a} and R^{1b} is not hydrogen;

R^2 is selected from the group consisting of: hydrogen,

haloalkyl,

C_1-C_6 alkyl,

C_1-C_6 alkyl-C_1-C_6 alkoxy,

C_1-C_6 alkyl-aryl,

C_2-C_6 alkenyl,

C_2-C_6 alkynyl,

$(CH_2)_n$-C_3-C_6 cycloalkyl,

C_1-C_6 alkoxy,

aryl, and ~~or~~

R^1 and R^2 together being a 5- to 8-membered heterocyclyl ring; ~~, and~~

wherein alkyl, aryl, alkenyl, alkynyl, cycloalkyl and alkoxy being optionally substituted with one or more groups independently selected from R^{15};

R^{2a} is selected from the group consisting of: hydrogen, halo, ~~or~~ C_1-C_6 alkyl, and ~~and~~ wherein R^2 and R^{2a} together being a 3- to 8-membered ring; ~~and~~ wherein alkyl being optionally substituted with one or more groups independently selected from R^{15};

R^3 is selected from the group consisting of: hydrogen,

halo,

cyano,

haloalkyl,

C_1–C_6 alkyl,

$(CH_2)_n$-C_3-C_6 cycloalkyl,

(C_1–C_4 alkyl)-heterocyclyl, and ~~wherein the heterocyclyl being optionally substituted with oxo,~~

(C_1–C_4 alkyl)-$NR^7C(O)_pR^9$; ~~, and~~

wherein alkyl, cycloalkyl and heterocyclyl being optionally substituted with one or more groups independently selected from R^{15};

R^4 and R^5 are each independently selected from the group consisting of:

hydrogen,

halo,

C_1-C_6 alkyl

C_1-C_6 alkoxy;

aryloxy;

$N(R^8)_2$,

SR^8 and~~or~~

R^4 and R^5 together being a 3- to 8-membered ring;

R^6 is selected from the group consisting of: hydrogen, C_1-C_6 alkyl and~~or~~ aminoalkyl;

R^{6A} is selected from the group consisting of: carboxamide, C_1-C_3 alkylnitrile, sulfonamide, acylsulfonamide and~~or~~ tetrazole;

R^7 is selected from the group consisting of: hydrogen and~~or~~ C_1-C_6 alkyl;

R^8 and R^9 are each independently selected from the group consisting of:

hydrogen, C_1-C_6 alkyl, aryl, heteroaryl, and~~or~~ heterocyclyl~~;, and~~

wherein aryl, heteroaryl and heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl and C_1-C_6 alkoxy;

R^{15} is selected from the group consisting of: hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl, C_1-C_6 alkoxy, $(CH_2)_n$-C_3-C_6 cycloalkyl, $N(R^8)_2$, $NR^8S(O)_2R^9$, $NR^8C(O)_pR^9$, $C(O)NR^8R^9$, $C(O)_pR^8$, SR^8, $S(O)_pR^8$ and~~or~~ $S(O)_2NR^8R^9$.

8. (Currently Amended) The compound of Claim 7, wherein X and Y are respectively S and O; S and CH_2; or CH_2 and O.

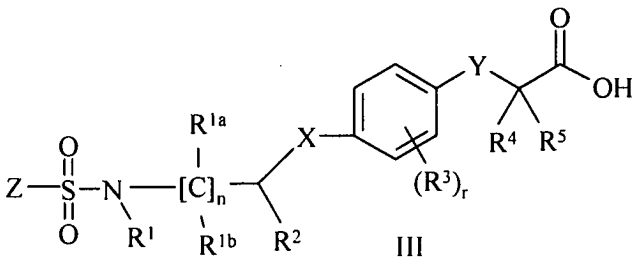
9. (Canceled)

10. (Canceled)

11. (Currently Amended) The compound of Claim 8~~Claim 10~~, wherein R^1 is C_3-C_6 alkyl or $(CH_2)_n$-C_3-C_6 cycloalkyl; R^2 and R^3 are each independently C_1-C_3 alkyl; and r is 1.

12. (Original) The compound Claim 11, wherein X is positioned para to Y; and R^3 is positioned ortho to Y.

13. (Currently Amended) The compound of Claim 7, wherein the compound having a structural Formula III,



III

orand pharmaceutically acceptable salts~~, solvates, hydrates or~~or stereoisomers thereof, wherein:

n is: 1 or 2;

r is: 1, 2, 3, or 4;

X is: S or CH_2;

Y is: CH_2 or O;

Z is: ~~aryl or a 5- to 10-membered heteroaryl,~~ benzothiophene;

wherein ~~aryl and heteroaryl being~~the benzothiophene is optionally substituted with one or more groups independently selected from R^{15};

R^1 and R^2 are each independently selected from the group consisting of~~:~~ C_1-C_6 alkyl and~~or~~ $(CH_2)_n$-C_3-C_6 cycloalkyl; and

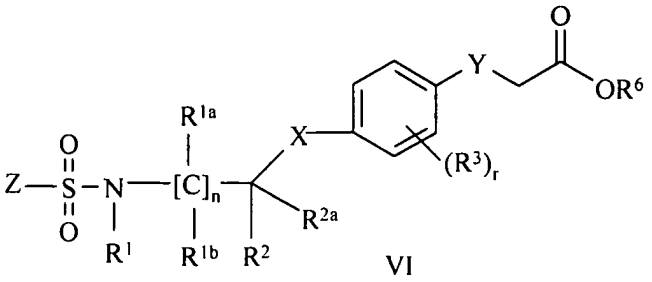
R^{1a} and R^{1b}, R^3, R^4 and R^5 are each independently selected from the group consisting of~~:~~ hydrogen and~~or~~ C_1-C_6 alkyl.

14. (Canceled)

15. (Canceled)

16. (Cancelled)

17. (Currently Amended) A compound having a structural Formula VI,



VI

orand pharmaceutically acceptable salts~~, solvates, hydrates~~ or stereoisomers thereof, wherein:

X is selected from the group consisting of~~:~~ a bond, CH_2, O, S and~~or~~ $S[O]_p$;

Y is selected from the group consisting of~~:~~ a bond, S, CH_2 and~~or~~ O;

Z is benzothiophene~~: heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S, and~~ wherein ~~heteroaryl being optionally~~the benzotihophene is optionally substituted with one or more groups selected from R^{15};

n is: 0, 1, 2 or 3;

n' is: 0, 1, 2, 3 or 4;

p is: 1 or 2;

r is: 1, 2, 3 or 4;

R^1 is selected from the group consisting of~~:~~ hydrogen,

haloalkyl,

C_1-C_6 alkyl,

C_1-C_6 alkyl-C_1-C_6 alkoxy,

C_1-C_6 alkyl-aryl,

C_2-C_6 alkenyl,

C_2-C_6 alkynyl,

$(CH_2)_n$-C_3-C_6 cycloalkyl,

C_1-C_6 alkoxy,

aryl, and~~or~~

R^1 and R^2 together being a 5- to 8-membered heterocyclyl ring~~; and~~

wherein alkyl, aryl, alkenyl, alkynyl, cycloalkyl and alkoxy being optionally substituted with one or more groups independently selected from R^{15};

R^{1a} and R^{1b} are each independently selected from the group consisting of: hydrogen, C_1-C_6 alkyl, and or R^1 and R^{1a}, R^1 and R^{1b}, R^2 and R^{1a}, R^2 and R^{1b} or R^{1a} and R^{1b} together being a 3- to 6-membered heterocyclyl or carbocyclyl ring; wherein where at least one of R^{1a} and R^{1b} is not hydrogen;

R^2 is selected from the group consisting of: hydrogen, haloalkyl, C_1-C_6 alkyl, C_1-C_6 alkyl-C_1-C_6 alkoxy, C_1-C_6 alkyl-aryl, C_2-C_6 alkenyl, C_2-C_6 alkynyl, $(CH_2)_n$-C_3-C_6 cycloalkyl, C_1-C_6 alkoxy, aryl, and or R^1 and R^2 together being a 5- to 8-membered heterocyclyl ring; and wherein alkyl, aryl, alkenyl, alkynyl, cycloalkyl and alkoxy being optionally substituted with one or more groups independently selected from R^{15};

R^{2a} is selected from the group consisting of: hydrogen, halo and or C_1-C_6 alkyl; and wherein R^2 and R^{2a} together being a 3- to 8-membered ring; and wherein alkyl being optionally substituted with one or more groups independently selected from R^{15};

R^3 is selected from the group consisting of: hydrogen, halo, cyano, haloalkyl, C_1-C_6 alkyl, $(CH_2)_n$-C_3-C_6 cycloalkyl,

(C_1–C_4 alkyl)-heterocyclyl, and~~wherein the heterocyclyl being optionally substituted with oxo;~~

(C_1–C_4 alkyl)-$NR^7C(O)_pR^9$,~~; and~~

wherein alkyl, cycloalkyl and heterocyclyl being optionally substituted with one or more groups independently selected from R^{15};

R^6 is selected from the group consisting of~~:~~ hydrogen, C_1-C_6 alkyl and~~or~~ aminoalkyl;

R^7 is selected from the group consisting of~~:~~ hydrogen and~~or~~ C_1-C_6 alkyl;

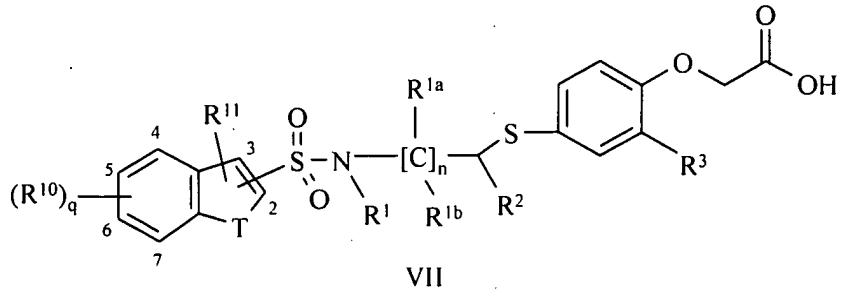
R^8 and R^9 are each independently selected from the group consisting of~~:~~

hydrogen, C_1-C_6 alkyl, aryl, heteroaryl, and~~or~~ heterocyclyl~~; and~~

wherein aryl, heteroaryl and heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl and C_1-C_6 alkoxy; and

R^{15} is selected from the group consisting of~~:~~ hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl, C_1-C_6 alkoxy, $N(R^8)_2$, $NR^8S(O)_2R^9$, $NR^8C(O)_pR^9$, $C(O)NR^8R^9$, $C(O)_pR^8$, SR^8, $S(O)_pR^8$ and~~or~~ $S(O)_2NR^8R^9$.

18. (Currently Amended) The compound of Claim17, wherein the compound having a structural Formula VII,



VII

or~~and~~ pharmaceutically acceptable salts or~~, solvates, hydrates or~~ stereoisomers thereof, wherein:

q is: 1, 2, 3, or 4;

T isis: O, NR^{1e} or S;

R^{1e} is: hydrogen or C_1-C_6 alkyl;

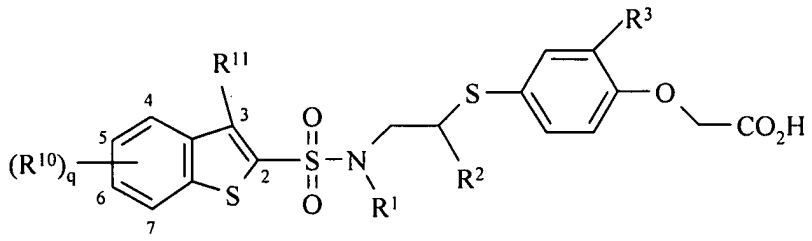
R^{10} and R^{11} are each independentlyselected from the group consisting of:

hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy,

C_1-C_6 alkyl andor C_1-C_6 alkoxy;; and

wherein alkyl, aryloxy, and alkoxy being optionally substituted with one or more groups independently selected from R^{15}.

19. (Currently Amended) The compound of Claim 18, wherein the compound having a structural Formula VIII,



VIII

orand pharmaceutically acceptable salts, solvates, hydrates or stereoisomers thereof, wherein:

q is: 1 or 2;

R^1 is selected from the group consisting of: C_3-C_5 alky andor $(CH_2)_n$-C_3-C_6 cycloalkyl;

R^2 and R^3 are each independently: C_1-C_3 alkyl;

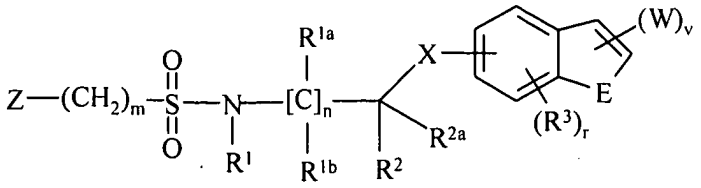
R^{10} is selected from the group consisting of :halo, haloalkyl and or C_1-C_3 alkyl;, and

wherein R^{10} being substituted at a position 5, or 6, or both 5 and 6 of benzothiophenyl ring; and

R^{11} is selected from the group consisting of: hydrogen andor C_1-C_6 alkyl.

20. (Original) The compound of Claim 19, wherein R^{10} is Cl, F, Br, CH_3 or CF_3 being substituted at a position 5 of benzothiophenyl ring.

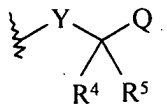
21. (Currently Amended) A compound having a structural Formula IX,



IX

orand pharmaceutically acceptable salts, ~~solvates, hydrates~~ or stereoisomers thereof, wherein:

E is selected from the group consisting of: O, S and~~or~~ NR^{14};

W is selected from the group consisting of:  , hydrogen, C_1-C_6 alkyl, $(CH_2)_n$-C_3-C_6 cycloalkyl, haloalkyl and~~or~~ acyl;

Q is selected from the group consisting of: -C(O)OR^6 and~~or~~ R^{6A};

X is selected from the group consisting of: a bond, C, O, S and~~or~~ $S[O]_p$;

Y is selected from the group consisting of: a bond, S, CH_2 ~~or~~and O;

Z is benzothiopene; and: ~~a) aliphatic group,~~

~~b) aryl,~~

~~c) a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S.~~

~~d) bi-aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl,~~

~~e) bi-heteroaryl, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and~~

~~f) heterocyclyl;~~

wherein ~~aliphatic group, aryl, heteroaryl, bi-aryl, bi-heteroaryl and heterocyclyl being~~the benzothiophene is optionally substituted with one or more groups independently selected from R^{15};

m and n' are each independently: 0, 1, 2, 3 or 4;

n is: 0, 1, 2 or 3;

p is: 1 or 2;

r is: 1, 2, 3 or 4;

v is: 1 or 2;

R^1 is selected from the group consisting of: hydrogen,

haloalkyl,

C_1-C_6 alkyl,

C_1-C_6 alkyl-C_1-C_6 alkoxy,

C_1-C_6 alkyl-aryl,

C_2-C_6 alkenyl,

C_2-C_6 alkynyl,

$(CH_2)_n$-C_3-C_6 cycloalkyl,

C_1-C_6 alkoxy,

aryl, and

R^1 and R^2 together being a 5- to 8-membered heterocyclyl ring;

wherein alkyl, aryl, alkenyl, alkynyl, cycloalkyl and alkoxy being optionally substituted with one or more groups independently selected from R^{15};

R^{1a} and R^{1b} are each independently selected from the group consisting of:

hydrogen,

C_1-C_6 alkyl, and

R^1 and R^{1a}, R^1 and R^{1b}, R^2 and R^{1a}, R^2 and R^{1b} or R^{1a} and R^{1b} together being a 3- to 6-membered heterocyclyl or carbocyclyl ring wherein at least one of R^{1a} and R^{1b} is not hydrogen;

R^2 is selected from the group consisting of: hydrogen,

haloalkyl,

C_1-C_6 alkyl,

C_1-C_6 alkyl-C_1-C_6 alkoxy,

C_1-C_6 alkyl-aryl,

C_2-C_6 alkenyl,

C_2-C_6 alkynyl,

$(CH_2)_n$-C_3-C_6 cycloalkyl,

C_1-C_6 alkoxy,

aryl, and~~or~~

R^1 and R^2 together being a 5- to 8-membered heterocyclyl ring;~~, and~~

wherein alkyl, aryl, alkenyl, alkynyl, cycloalkyl and alkoxy being optionally substituted with one or more groups independently selected from R^{15};

R^{2a} is selected from the group consisting of~~:~~ hydrogen, halo,~~ or~~ C_1-C_6 alkyl and~~ and~~ wherein R^2 and R^{2a} together being a 3- to 8-membered ring;~~ and~~ wherein alkyl being optionally substituted with one or more groups independently selected from R^{15};

R^3 is selected from the group consisting of~~:~~ hydrogen,

halo,

cyano,

haloalkyl,

C_1–C_6 alkyl,

$(CH_2)_n$-C_3-C_6 cycloalkyl,

(C_1–C_4 alkyl)-heterocyclyl,~~ wherein the heterocyclyl being optionally substituted with oxo,~~ and

(C_1–C_4 alkyl)-$NR^7C(O)_pR^9$;~~ and~~

wherein alkyl, cycloalkyl and heterocyclyl being optionally substituted with one or more groups independently selected from R^{15};

R^4 and R^5 are each independently selected from the group consisting of~~:~~

hydrogen,

halo,

C_1-C_6 alkyl

C_1-C_6 alkoxy;

aryloxy;

$N(R^8)_2$,

SR^8 and~~or~~

R^4 and R^5 together being a 3- to 8-membered ring;

R^6 is selected from the group consisting of~~:~~ hydrogen, C_1-C_6 alkyl and~~or~~ aminoalkyl;

R^{6A} is selected from the group consisting of: carboxamide, C_1-C_3 alkylnitrile, sulfonamide, acylsulfonamide and tetrazole;

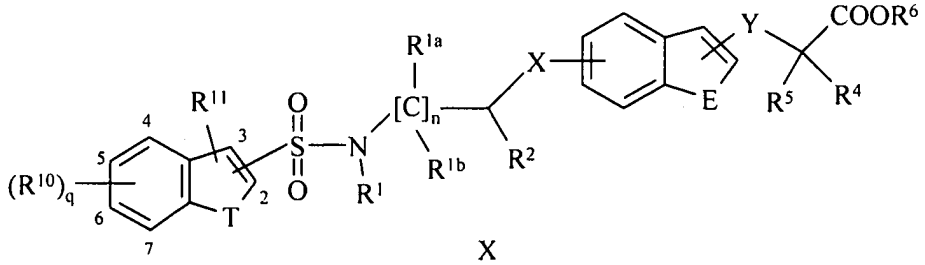
R^7 is selected from the group consisting of: hydrogen and C_1-C_6 alkyl;

R^8 and R^9 are each independently selected from the group consisting of: hydrogen, C_1-C_6 alkyl, aryl, heteroaryl, and heterocyclyl; and wherein aryl, heteroaryl and heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl and C_1-C_6 alkoxy;

R^{14} is selected from the group consisting of: hydrogen, aryl, C_1-C_6 alkyl, and C_1-C_6 alkyl-$COOR^6$; and wherein aryl and alkyl being optionally substituted with one or more groups independently selected from R^{15}; and

R^{15} is selected from the group consisting of: hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1-C_6 alkyl, C_1-C_6 alkoxy, $(CH_2)_n$-C_3-C_6 cycloalkyl, $N(R^8)_2$, $NR^8S(O)_2R^9$, $NR^8C(O)_pR^9$, $C(O)NR^8R^9$, $C(O)_pR^8$, SR^8, $S(O)_pR^8$ and $S(O)_2NR^8R^9$.

22. (Currently Amended) The compound of Claim 21, wherein the compound having a structural Formula X:



X

and pharmaceutically acceptable salts, solvates, hydrates or stereoisomers thereof, wherein:

n and q are each independently: 1, 2, 3 or 4;

T is: O, NR^{1e} or S;

X is selected from the group consisting of: CH_2, O and~~or~~ S;

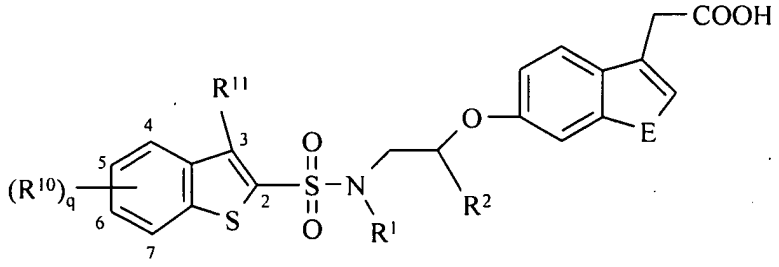
R^1 is selected from the group consisting of: hydrogen, C_1-C_6 alkyl and~~or~~ $(CH_2)_n$-C_3-C_6 cycloalkyl;

R^{1a}, R^{1b}, R^{1c} and R^2 are each independently selected from the group consisting of: hydrogen and~~or~~ C_1-C_6 alkyl; and

R^{10} and R^{11} are each independently selected from the group consisting of:

hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, C_1-C_6 alkyl and~~or~~ C_1-C_6 alkoxy; ~~and~~ wherein alkyl, alkoxy and aryloxy being optionally substituted with one or more groups selected from R^{15}.

23. (Currently Amended) The compound of Claim 22, wherein the compound having a structural Formula XI:



XI

or~~and~~ pharmaceutically acceptable salts~~, solvates, hydrates~~ or stereoisomers thereof, wherein:

q is 1 or 2;

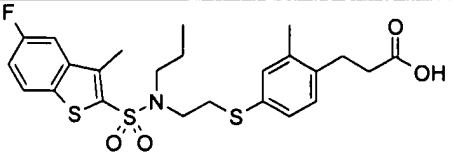
E is selected from the group consisting of O, S and~~or~~ NR^{14};

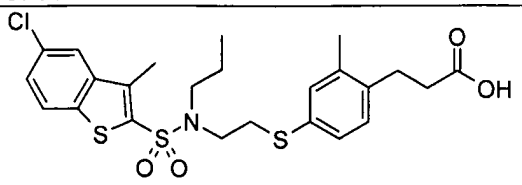
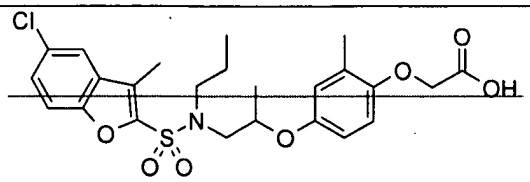
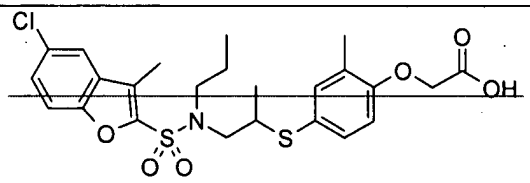
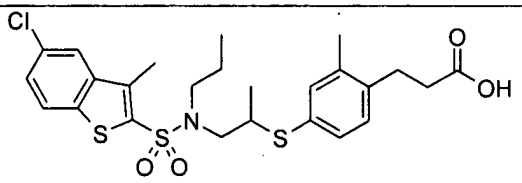
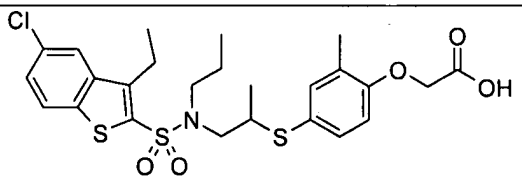
R^1, R^2 and R^{11} are each independently selected from the group consisting of: C_1-C_4 alkyl;

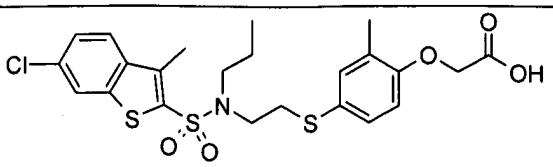
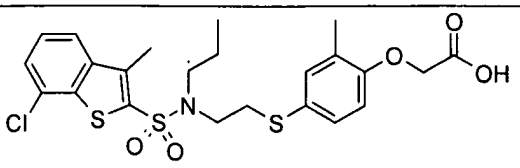
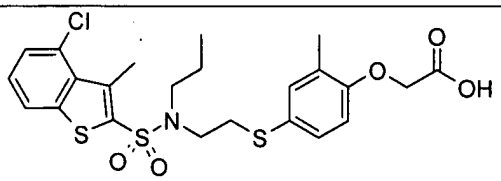
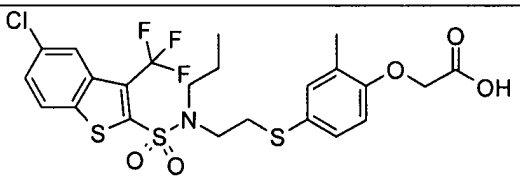
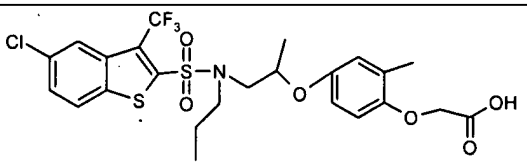
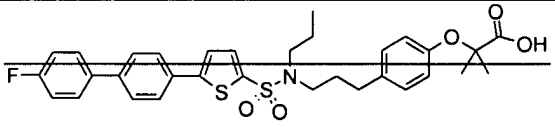
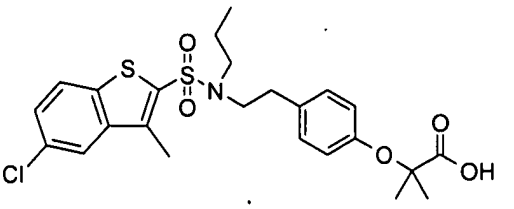
R^{10} is selected from the group consisting of: Cl, F, Br, CH_3 and~~or~~ CF_3; ~~and~~ wherein R^{10} being substituted at a position 5, or 6, or both 5 and 6 of benzothiophenyl ring; and

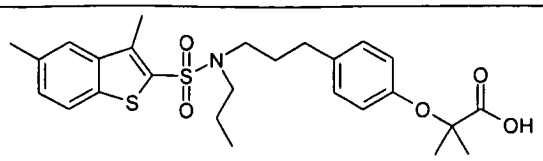
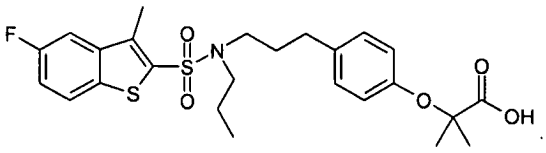
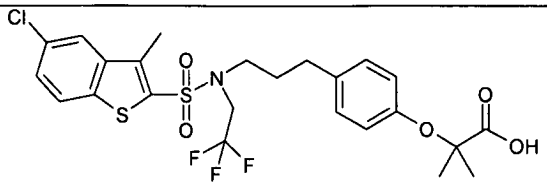
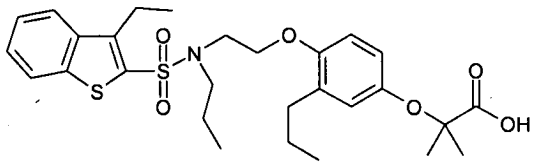
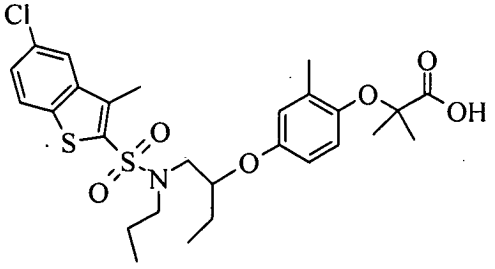
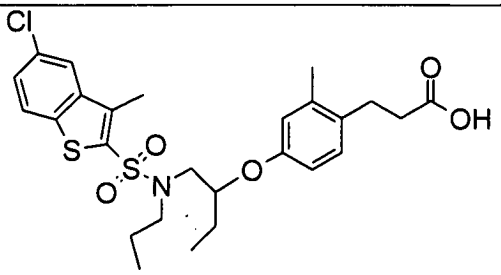
R^{14} is selected from the group consisting of: hydrogen, C_1-C_6 alkyl and~~or~~ aryl.

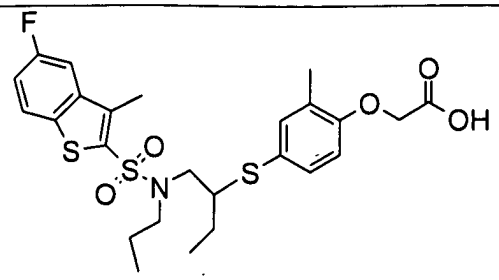
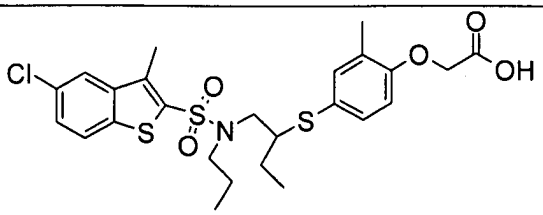
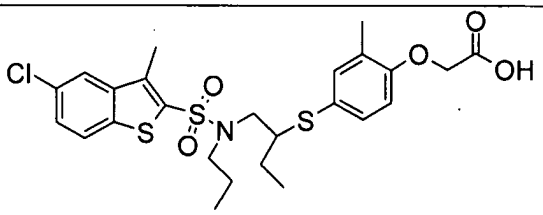
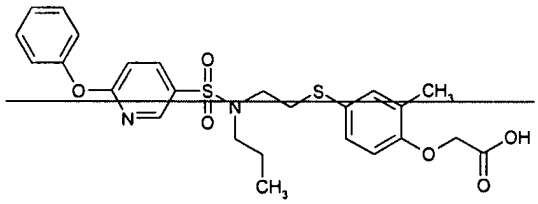
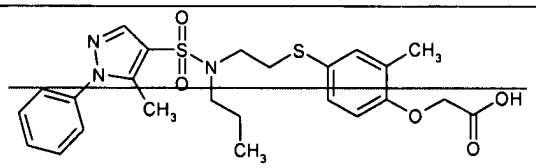
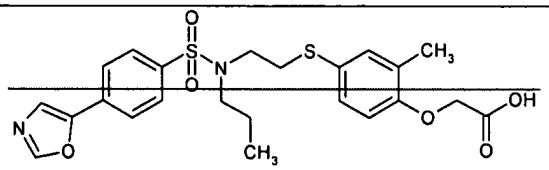
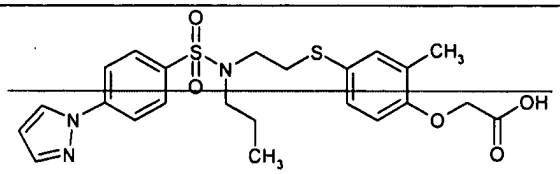
24. (Currently Amended) A compound selected from the group consisting of ~~No. 1-120 and 121:~~

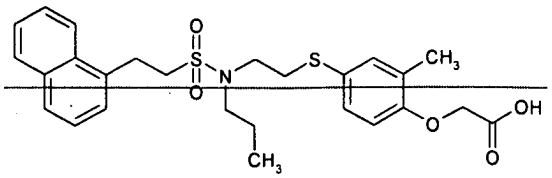
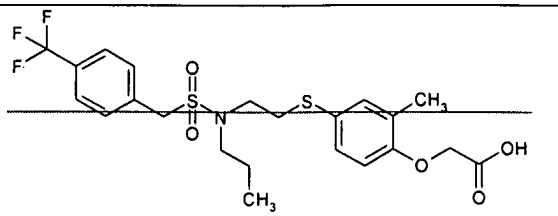
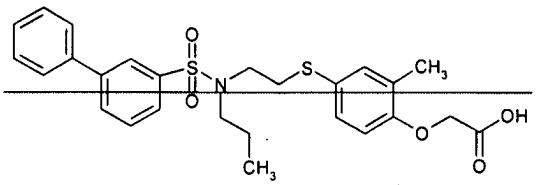
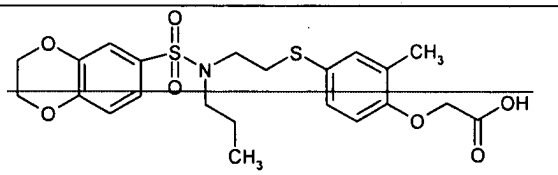
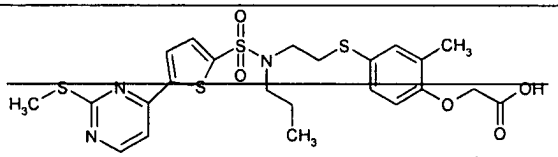
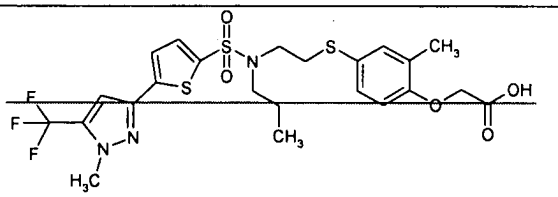
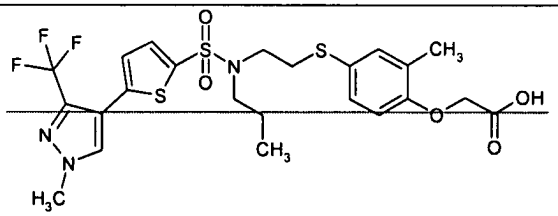
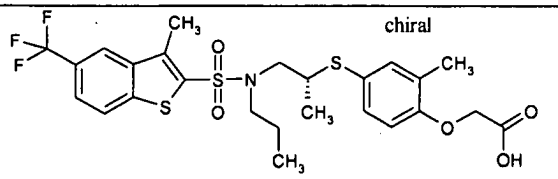
No.	Structure	Name
1		3-(4-{2-[(5-Fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenyl)-propionic acid

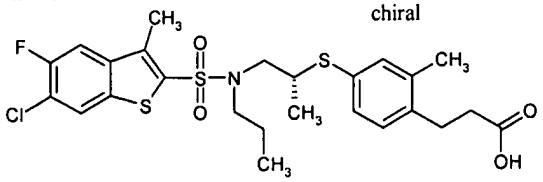
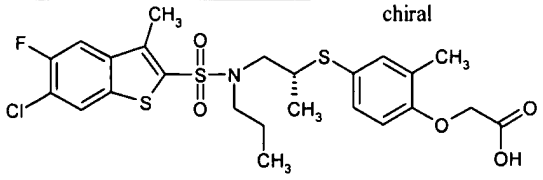
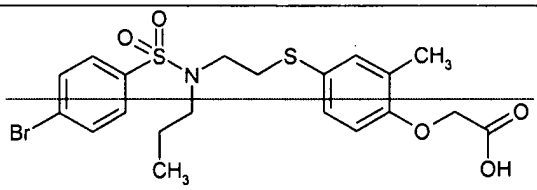
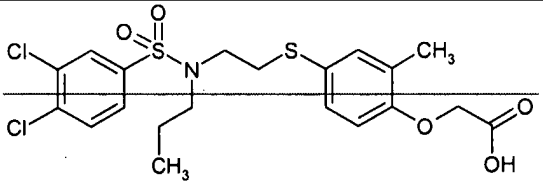
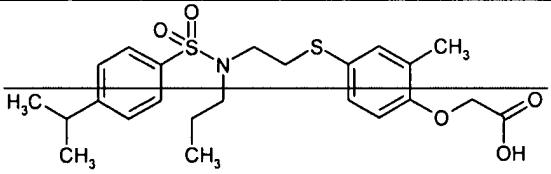
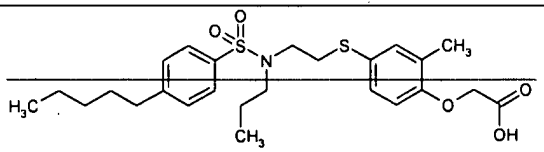
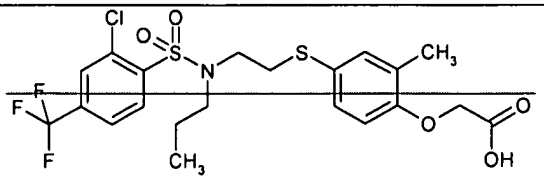
No.	Structure	Name
2		3-(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenyl)-propionic acid
~~3~~		~~(4-{2-[(5-Chloro-3-methyl-benzofuran-2-sulfonyl)-propyl-amino]-1-methyl-ethoxy}-2-methyl-phenoxy)-acetic acid~~
4		~~(4-{2-[(5-Chloro-3-methyl-benzofuran-2-sulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
5		3-(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenyl)-propionic acid
6		(4-{2-[(5-Chloro-3-ethyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid

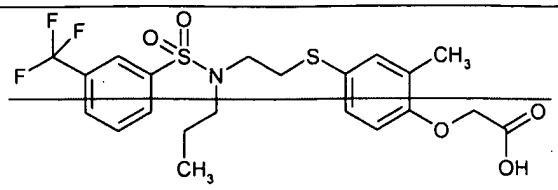
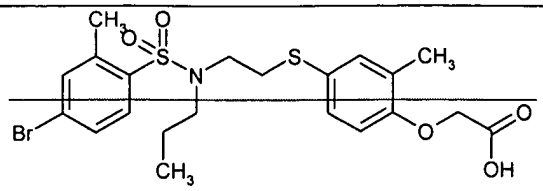
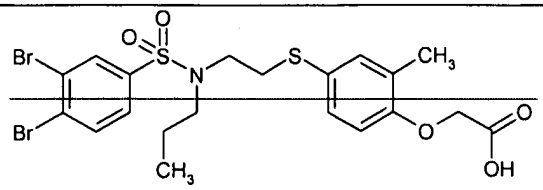
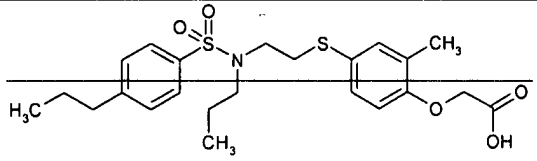
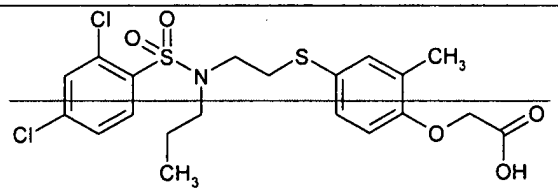
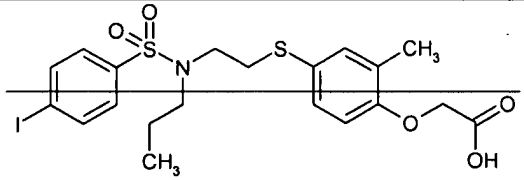
7		4-{2-[(6-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
8		4-{2-[(7-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
9		(4-{2-[(4-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
10		(4-{2-[(5-Chloro-3-trifluoromethyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
11		(4-{2-[(5-Chloro-3-trifluoromethyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethoxy}-2-methyl-phenoxy)-acetic acid
~~12~~		~~2-[4-(3-{[5-(4'-Fluoro-biphenyl-4-yl)-thiophene-2-sulfonyl]-propyl-amino}-propyl)-phenoxy]-2-methyl-propionic acid~~
13		2-(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethyl}-phenoxy)-2-methyl-propionic acid

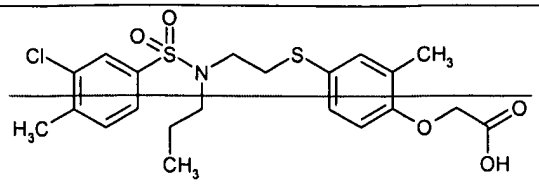
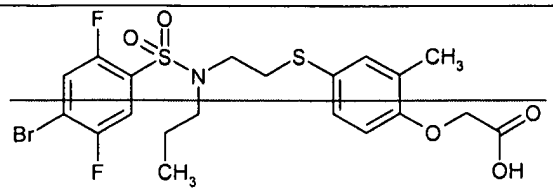
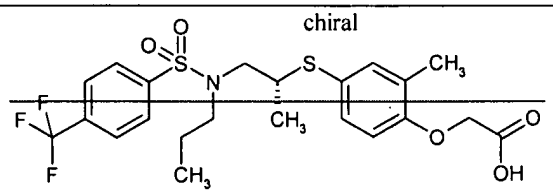
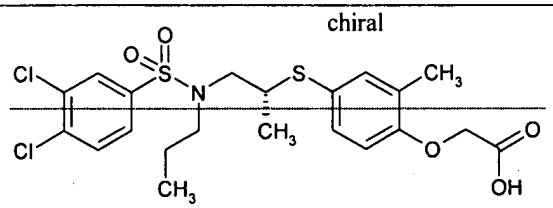
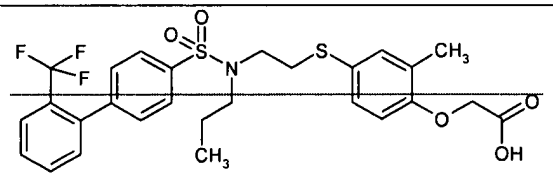
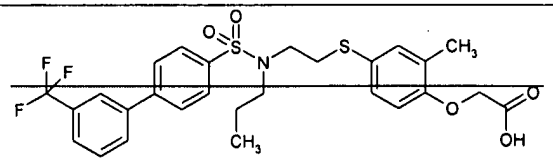
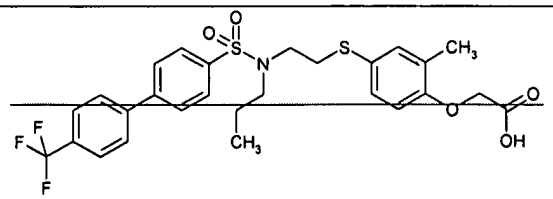
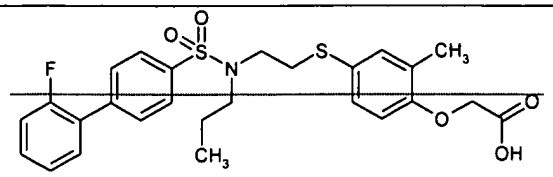
14		2-(4-{3-[(3,5-Dimethyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-propyl}-phenoxy)-2-methyl-propionic acid
15		2-(4-{3-[(5-Fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-propyl}-phenoxy)-2-methyl-propionic acid
16		2-(4-{3-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-(2,2,2-trifluoro-ethyl)-amino]-propyl}-phenoxy)-2-methyl-propionic acid
17		2-(4-{2-[(3-Ethyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethoxy}-3-propyl-phenoxy)-2-methyl-propionic acid
18		2-[4-(1-{[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-methyl}-propoxy)-2-methyl-phenoxy]-2-methyl-propionic acid
19		3-[4-(1-{[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-methyl}-propoxy)-2-methyl-phenyl]-propionic acid

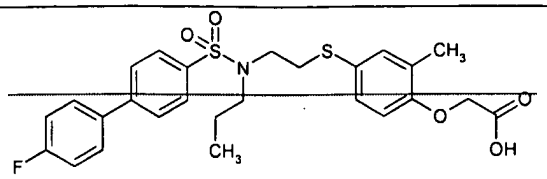
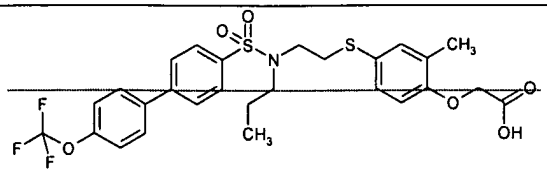
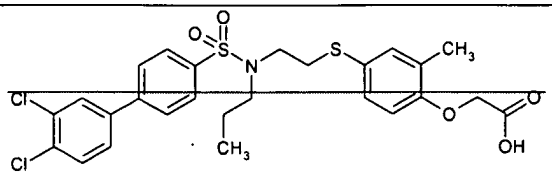
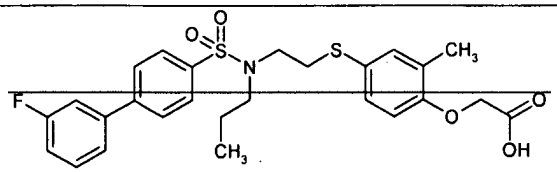
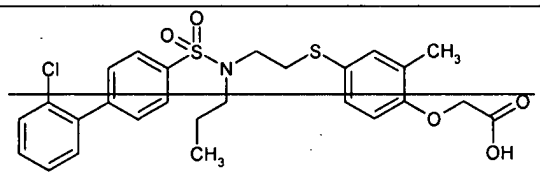
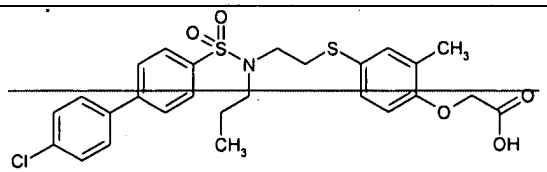
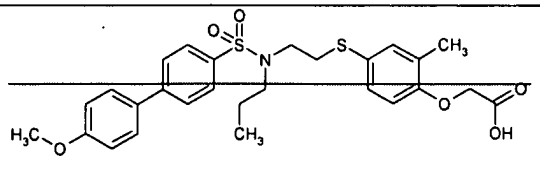
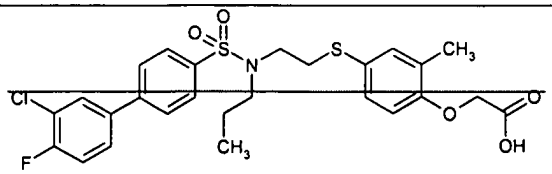
20		[4-(1-{[(5-Fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-methyl}-propylsulfanyl)-2-methyl-phenoxy]-acetic acid
21		[4-(1-{[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-methyl}-propylsulfanyl)-2-methyl-phenoxy]-acetic acid
22		[4-(1-{[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-methyl}-propylsulfanyl)-2-methyl-phenoxy]-acetic acid
~~23~~		~~(2-Methyl-4-{2-[(6-phenoxy-pyridine-3-sulfonyl)-propyl-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~24~~		~~(2-Methyl-4-{2-[(5-methyl-1-phenyl-1H-pyrazole-4-sulfonyl)-propyl-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~25~~		~~(2-Methyl-4-{2-[(4-oxazol-5-yl-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~26~~		~~(2-Methyl-4-{2-[propyl-(4-pyrazol-1-yl-benzenesulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~

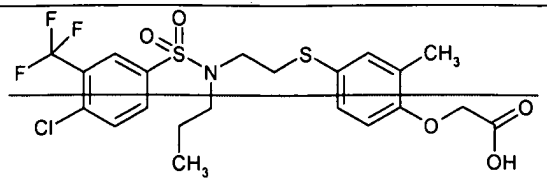
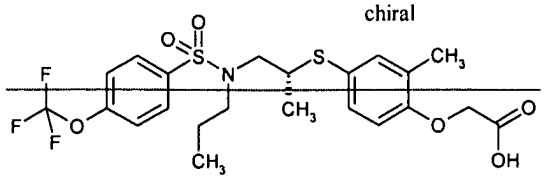
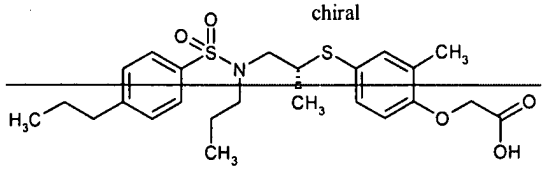
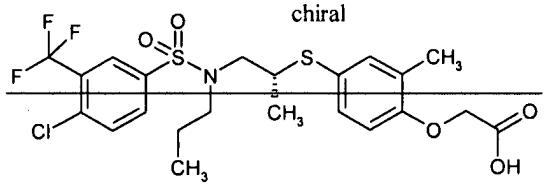
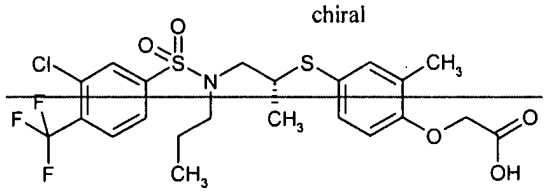
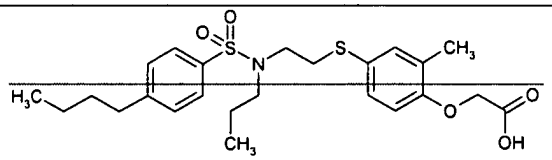
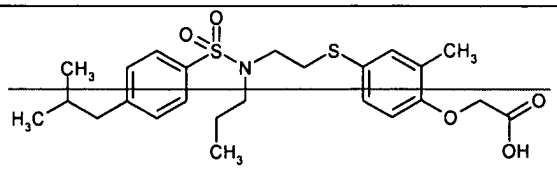
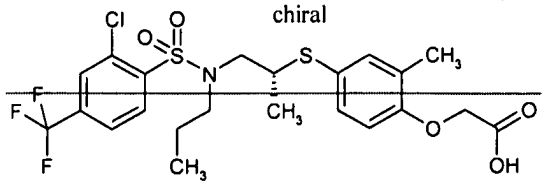
~~27~~		~~(2-Methyl-4-{2-[(2-naphthalen-1-yl-ethanesulfonyl)-propyl-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~28~~		~~(2-Methyl-4-{2-[propyl-(4-trifluoromethylphenylmethanesulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~29~~		~~(4-{2-[(Biphenyl-3-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~30~~		~~(4-{2-[(2,3-Dihydro-benzo[1,4]dioxine-6-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~31~~		~~[2-Methyl-4-(2-{[5-(2-methylsulfanyl-pyrimidin-4-yl)-thiophene-2-sulfonyl]-propyl-amino}-ethylsulfanyl)-phenoxy]-acetic acid~~
~~32~~		~~[2-Methyl-4-(2-{[5-(1-methyl-5-trifluoromethyl-1H-pyrazol-3-yl)-thiophene-2-sulfonyl]-propyl-amino}-ethylsulfanyl)-phenoxy]-acetic acid~~
~~33~~		~~[2-Methyl-4-(2-{[5-(1-methyl-3-trifluoromethyl-1H-pyrazol-4-yl)-thiophene-2-sulfonyl]-propyl-amino}-ethylsulfanyl)-phenoxy]-acetic acid~~
34		(R)-(2-Methyl-4-{1-methyl-2-[(3-methyl-5-trifluoromethyl-benzo[b]thiophene-2-sulfonyl)-propyl-

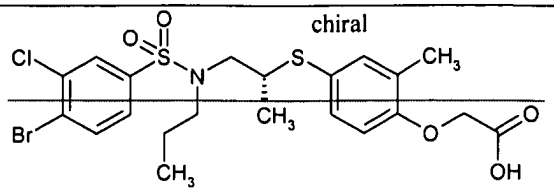
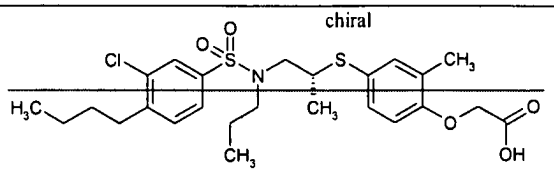
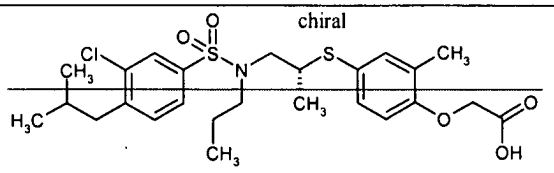
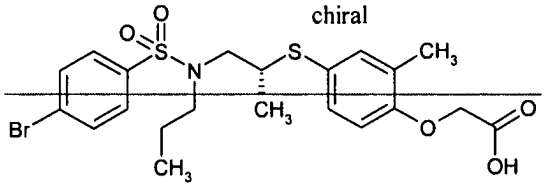
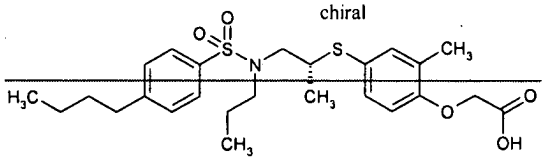
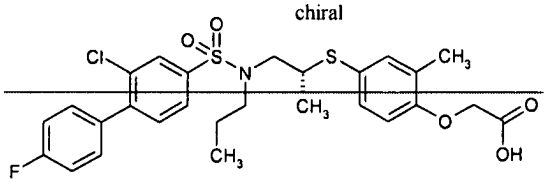
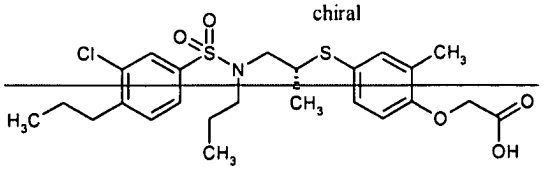
		amino]-ethylsulfanyl}-phenoxy)-acetic acid
35		(R)-3-(4-{2-[(6-Chloro-5-fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenyl)-propionic acid
36		(R)-(4-{2-[(6-Chloro-5-fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
~~37~~		~~(4-{2-[(4-Bromo-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~38~~		~~(4-{2-[(3,4-Dichloro-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~39~~		~~(4-{2-[(4-Isopropyl-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~40~~		~~(2-Methyl-4-{2-[(4-pentyl-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~41~~		~~(4-{2-[(2-Chloro-4-trifluoromethyl-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~

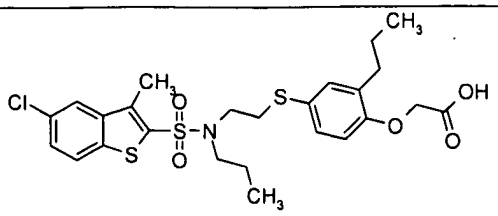
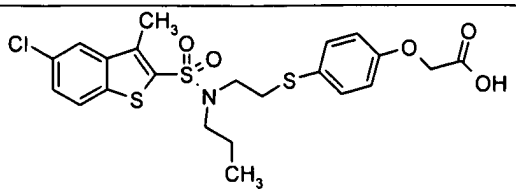
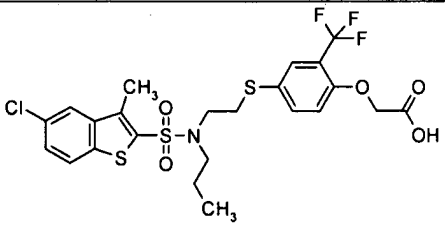
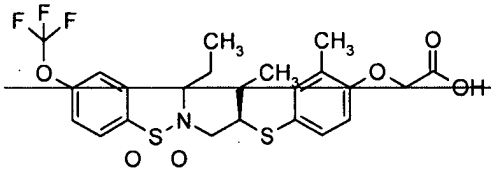
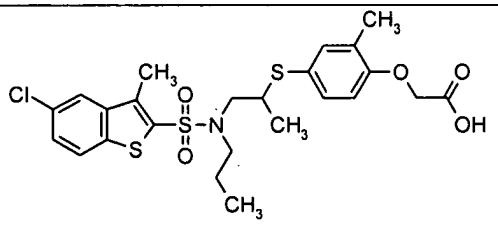
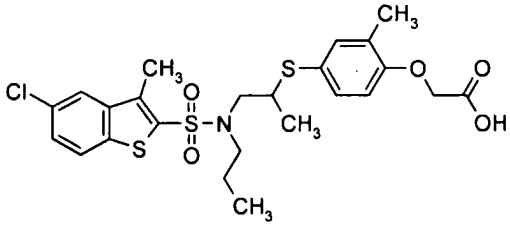
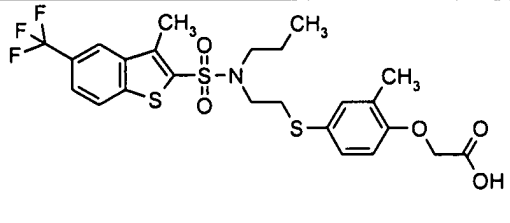
~~42~~		~~(2-Methyl-4-{2-[propyl-(3-trifluoromethyl-benzenesulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~43~~		~~(4-{2-[(4-Bromo-2-methyl-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~44~~		~~(4-{2-[(3,4-Dibromo-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~45~~		~~(2-Methyl-4-{2-[propyl-(4-propyl-benzenesulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~46~~		~~(4-{2-[(2,4-Dichloro-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~47~~		~~(4-{2-[(4-Iodo-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~

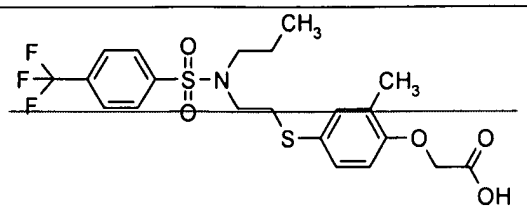
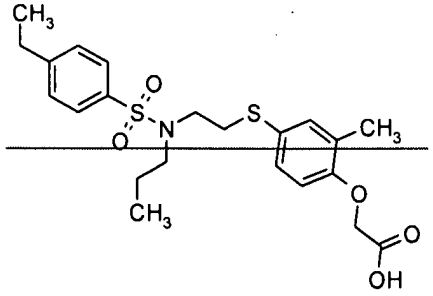
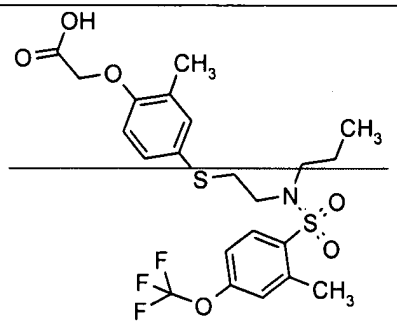
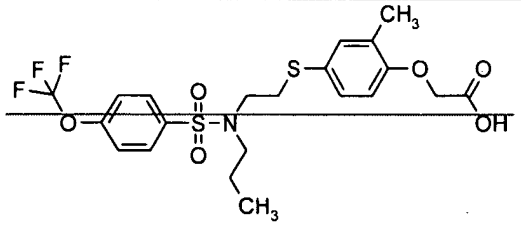
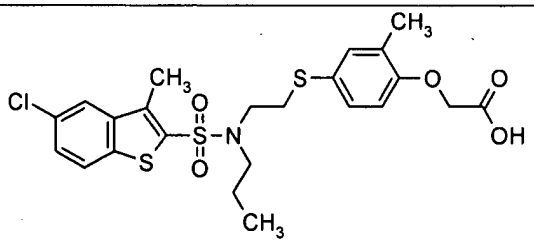
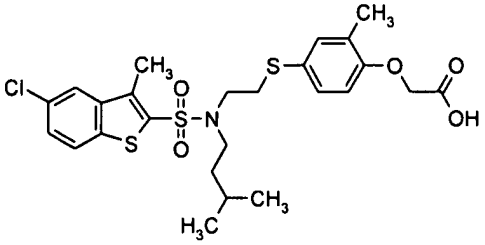
~~48~~		~~(4-{2-[(3-Chloro-4-methyl-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~49~~		~~(4-{2-[(4-Bromo-2,5-difluoro-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~50~~		~~(2-Methyl-4-{1-methyl-2-[propyl-(4-trifluoromethyl-benzenesulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~51~~		~~(4-{2-[(3,4-Dichloro-benzenesulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~52~~		~~(2-Methyl-4-{2-[propyl-(2'-trifluoromethyl-biphenyl-4-sulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~53~~		~~(2-Methyl-4-{2-[propyl-(3'-trifluoromethyl-biphenyl-4-sulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~54~~		~~(2-Methyl-4-{2-[propyl-(4'-trifluoromethyl-biphenyl-4-sulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~55~~		~~(4-{2-[(2'-Fluoro-biphenyl-4-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~

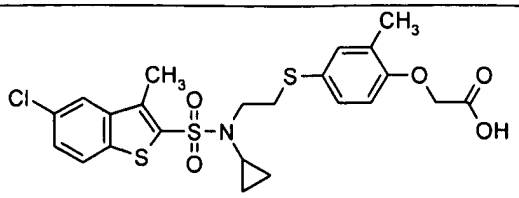
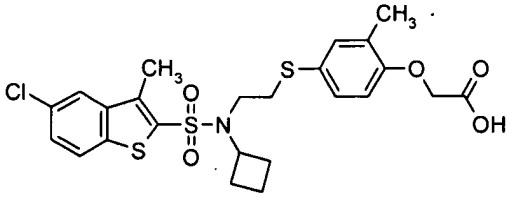
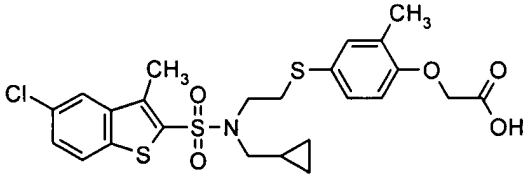
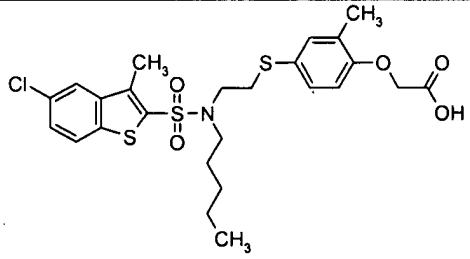
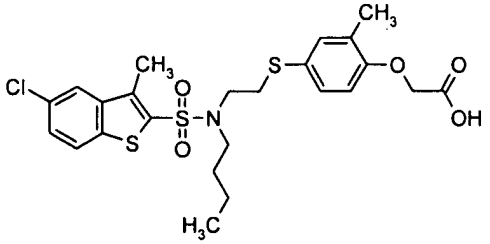
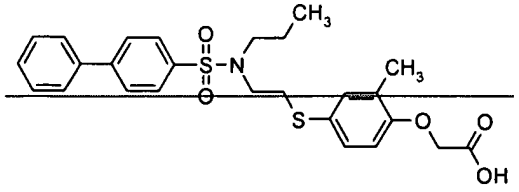
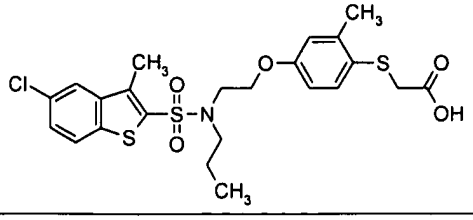
~~56~~		~~(4-{2-[(4'-Fluoro-biphenyl-4-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~57~~		~~(2-Methyl-4-{2-[propyl-(4'-trifluoromethoxy-biphenyl-4-sulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~58~~		~~(4-{2-[(3',4'-Dichloro-biphenyl-4-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~59~~		~~(4-{2-[(3'-Fluoro-biphenyl-4-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~60~~		~~(4-{2-[(2'-Chloro-biphenyl-4-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~61~~		~~(4-{2-[(4'-Methoxy-biphenyl-4-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~62~~		~~(4-{2-[(4'-Methoxy-biphenyl-4-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~63~~		~~(4-{2-[(3'-Chloro-4'-fluoro-biphenyl-4-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~

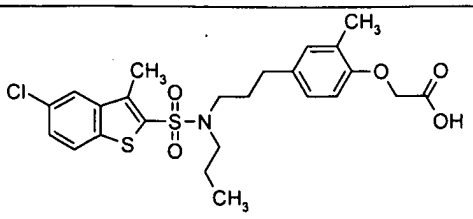
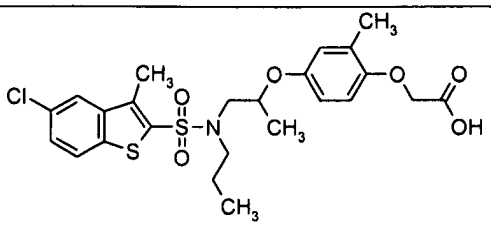
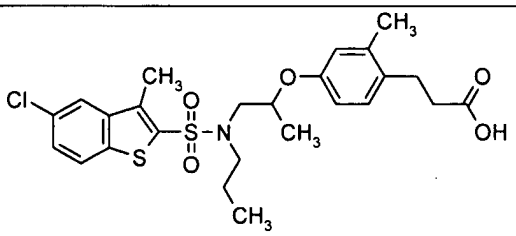
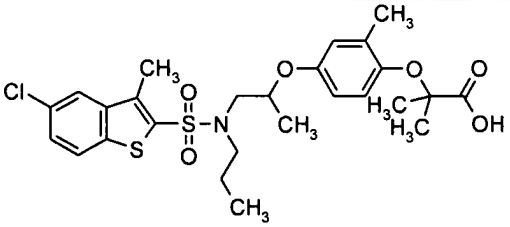
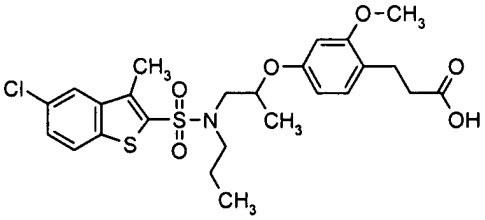
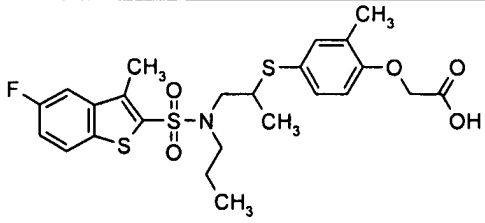
~~64~~		~~(4-{2-[(4-Chloro-3-trifluoromethyl-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~65~~		~~(2-Methyl-4-{1-methyl-2-[propyl-(4-trifluoromethoxy-benzenesulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~66~~		~~(2-Methyl-4-{1-methyl-2-[propyl-(4-propyl-benzenesulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~67~~		~~(4-{2-[(4-Chloro-3-trifluoromethyl-benzenesulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~68~~		~~(4-{2-[(3-Chloro-4-trifluoromethyl-benzenesulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~69~~		~~(4-{2-[(4-Butyl-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~70~~		~~(4-{2-[(4-Isobutyl-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~71~~		~~(4-{2-[(2-Chloro-4-trifluoromethyl-benzenesulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-~~

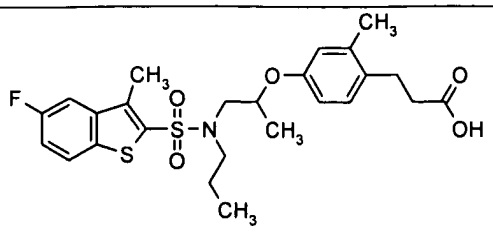
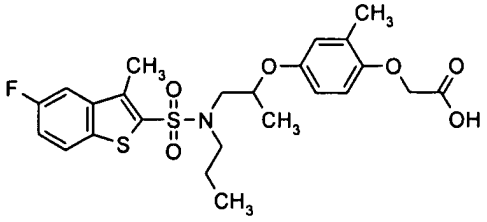
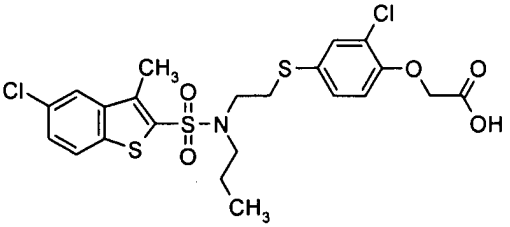
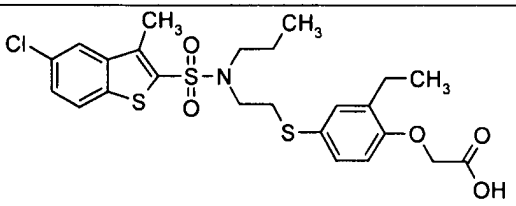
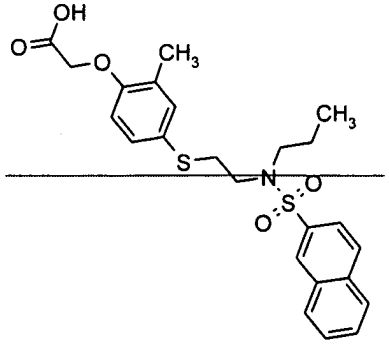
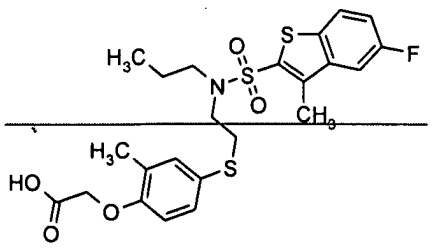
		~~2-methyl-phenoxy) acetic acid~~
~~72~~		~~(4-{2-[(4-Bromo-3-chloro-benzenesulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy) acetic acid~~
~~73~~		~~(4-{2-[(4-Butyl-3-chloro-benzenesulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy) acetic acid~~
~~74~~		~~(4-{2-[(3-Chloro-4-isobutyl-benzenesulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy) acetic acid~~
~~75~~		~~(4-{2-[(4-Bromo-benzenesulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy) acetic acid~~
~~76~~		~~(4-{2-[(4-Butyl-benzenesulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy) acetic acid~~
~~77~~		~~(4-{2-[(2-Chloro-4'-fluoro-biphenyl-4-sulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~78~~		~~(4-{2-[(3-Chloro-4-propyl-benzenesulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy) acetic acid~~

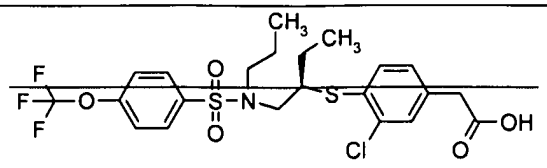
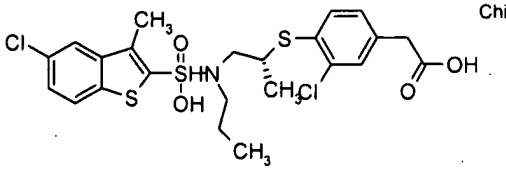
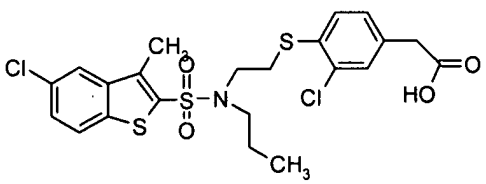
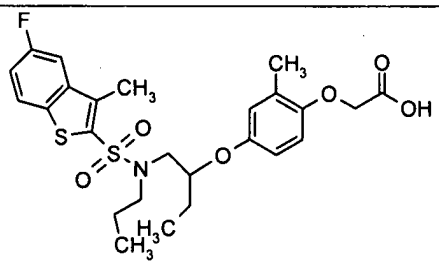
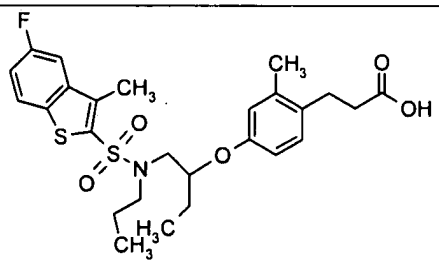
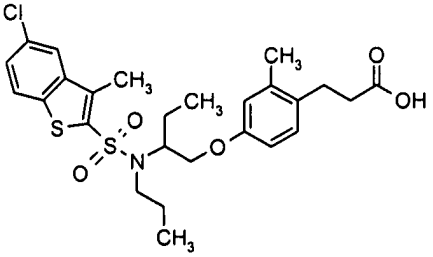
79		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-propyl-phenoxy)-acetic acid
80		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-phenoxy)-acetic acid
81		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-trifluoromethyl-phenoxy)-acetic acid
~~82~~		~~[2-Methyl-4-(1-{[propyl-(4-trifluoromethoxy-benzenesulfonyl)-amino]-methyl}-propylsulfanyl)-phenoxy]-acetic acid~~
83		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
84		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
85		(2-Methyl-4-{2-[(3-methyl-5-trifluoromethyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-phenoxy)-acetic acid

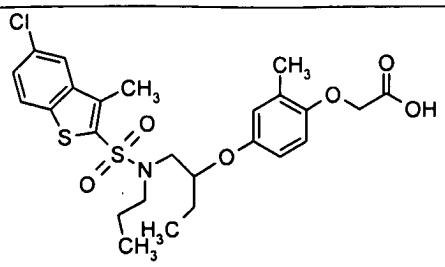
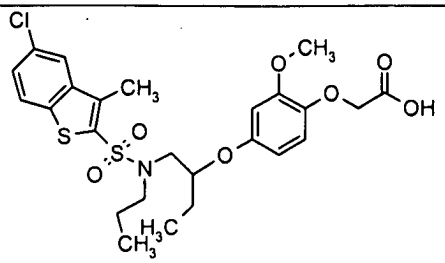
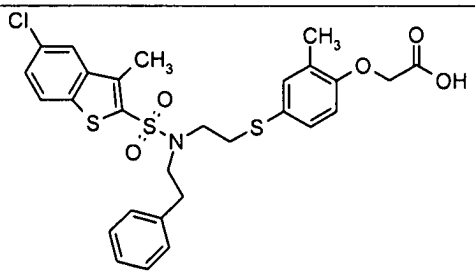
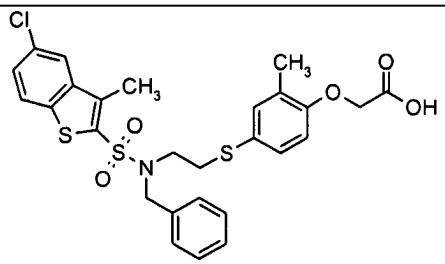
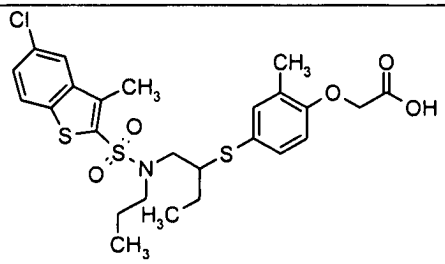
~~86~~		~~(2-Methyl-4-{2-[propyl-(4-trifluoromethyl-benzenesulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~87~~		~~(4-{2-[(4-Ethyl-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
~~88~~		~~(2-Methyl-4-{2-[(2-methyl-4-trifluoromethoxy-benzenesulfonyl)-propyl-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~89~~		~~(2-Methyl-4-{2-[propyl-(4-trifluoromethoxy-benzenesulfonyl)-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
90		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
91		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-(3-methyl-butyl)-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid

92		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-cyclopropyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
93		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-cyclobutyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
94		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-cyclopropylmethyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
95		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-pentyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
96		(4-{2-[Butyl-(5-chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
~~97~~		~~(4-{2-[(Biphenyl-4-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~
98		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethoxy}-2-methyl-phenylsulfanyl)-

		acetic acid
99		(4-{3-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-propyl}-2-methyl-phenoxy)-acetic acid
100		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethoxy}-2-methyl-phenoxy)-acetic acid
101		3-(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethoxy}-2-methyl-phenyl)-propionic acid
102		2-(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethoxy}-2-methyl-phenoxy)-2-methyl-propionic acid
103		3-(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethoxy}-2-methoxy-phenyl)-propionic acid
104		(4-{2-[(5-Fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid

105		3-(4-{2-[(5-Fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethoxy}-2-methyl-phenyl)-propionic acid
106		(4-{2-[(5-Fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethoxy}-2-methyl-phenoxy)-acetic acid
107		(2-Chloro-4-{2-[(5-chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-phenoxy)-acetic acid
108		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-ethyl-phenoxy)-acetic acid
~~109~~		~~(2-Methyl-4-{2-[(naphthalene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-phenoxy)-acetic acid~~
~~110~~		~~(4-{2-[(5-Fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid~~

~~111~~		~~[3-Chloro-4-(1-{[propyl-(4-trifluoromethoxy-benzenesulfonyl)-amino]-methyl}-propylsulfanyl)-phenyl]-acetic acid~~
112	 Chiral	(R)-(3-Chloro-4-{2-[(5-chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-1-methyl-ethylsulfanyl}-phenyl)-acetic acid
113		(3-Chloro-4-{2-[(5-chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-ethylsulfanyl}-phenyl)-acetic acid
114		[4-(1-{[(5-Fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-methyl}-propoxy)-2-methyl-phenoxy]-acetic acid
115		3-[4-(1-{[(5-Fluoro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-methyl}-propoxy)-2-methyl-phenyl]-propionic acid
116		3-(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-butoxy}-2-methyl-phenyl)-propionic acid

117		[4-(1-{[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-methyl}-propoxy)-2-methyl-phenoxy]-acetic acid
118		[4-(1-{[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-methyl}-propoxy)-2-methoxy-phenoxy]-acetic acid
119		(4-{2-[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-phenethyl-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid
120		(4-{2-[Benzyl-(5-chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-amino]-ethylsulfanyl}-2-methyl-phenoxy)-acetic acid; and
121		[4-(1-{[(5-Chloro-3-methyl-benzo[b]thiophene-2-sulfonyl)-propyl-amino]-methyl}-propylsulfanyl)-2-methyl-phenoxy]-acetic acid

25. (Currently Amended) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and at least one compound of Claim 1 or pharmaceutically acceptable salts~~, solvates or hydrates~~ thereof.

26. (Cancelled)

27. (Cancelled)

28. (Cancelled)

29. (Cancelled)

30. (Cancelled)

31. (Cancelled)

32. (Cancelled)

33. (Cancelled)

34. (Previously Presented) A method for lowering blood-glucose in a mammal comprising the step of administering an effective amount of at least one compound of Claim 1.

35. (Cancelled)

36. (Cancelled)

37. (Cancelled)

38. (Cancelled)

39. (Cancelled)

40. (Cancelled)